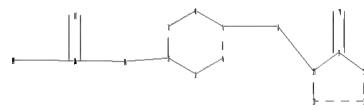
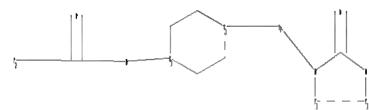


10/589515

=>

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chain nodes :

8 14 15 16 17 19

ring nodes :

1 2 3 4 5 6 9 10 11 12 13

chain bonds :

2-15 5-8 8-13 9-14 15-16 16-17 16-19

ring bonds :

1-2 1-6 2-3 3-4 4-5 5-6 9-10 9-13 10-11 11-12 12-13

exact/norm bonds :

1-2 1-6 2-3 2-15 3-4 4-5 5-6 5-8 8-13 9-10 9-13 9-14 10-11 11-12
12-13 15-16 16-17 16-19

G1:C,N

G2:O,N

Match level :

1:Atom 2:Atom 3:Atom 4:Atom 5:Atom 6:Atom 8:Atom 9:Atom 10:Atom 11:Atom
12:Atom 13:Atom 14:CLASS 15:CLASS 16:CLASS 17:CLASS 19:CLASS

Generic attributes :

8:

Saturation : Unsaturated
Number of Carbon Atoms : less than 7
Type of Ring System : Monocyclic

10/589515

L1 STRUCTURE UPLOADED

=> s 11
SAMPLE SEARCH INITIATED 13:21:17 FILE 'REGISTRY'
SAMPLE SCREEN SEARCH COMPLETED - 1040 TO ITERATE

100.0% PROCESSED 1040 ITERATIONS 0 ANSWERS
SEARCH TIME: 00.00.01

FULL FILE PROJECTIONS: ONLINE **COMPLETE**
BATCH **COMPLETE**
PROJECTED ITERATIONS: 18866 TO 22734
PROJECTED ANSWERS: 0 TO 0

L2 0 SEA SSS SAM L1

=> s 11 sss full
FULL SEARCH INITIATED 13:21:26 FILE 'REGISTRY'
FULL SCREEN SEARCH COMPLETED - 20152 TO ITERATE

100.0% PROCESSED 20152 ITERATIONS 2 ANSWERS
SEARCH TIME: 00.00.01

L3 2 SEA SSS FUL L1

=> file caplus
COST IN U.S. DOLLARS SINCE FILE TOTAL
SESSION
FULL ESTIMATED COST ENTRY 178.82 179.03

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FILE LAST UPDATED: 28 Apr 2008 (20080428/ED)

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=> s 13
L4 1 L3

10/589515

=> d 14

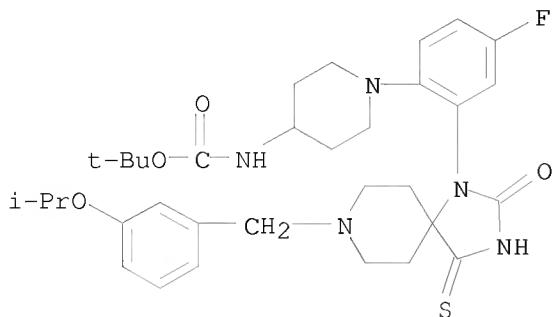
L4 ANSWER 1 OF 1 CAPLUS COPYRIGHT 2008 ACS on STN
AN 2006:383818 CAPLUS
DN 144:432808
TI Preparation of spiropiperidine compounds as β -secretase inhibitors for the treatment of Alzheimer's disease
IN Barrow, James C.; Coburn, Craig A.; Egbertson, Melissa S.; McGaughey, Georgia B.; McWherter, Melody A.; Neilson, Lou Anne; Selnick, Harold G.; Stauffer, Shaun R.; Yang, Zhi-Qiang; Yang, Wenjin; Lu, Wanli; Fahr, Bruce; Rittle, Kenneth E.
PA Merck & Co., Inc., USA; Sunesis Pharmaceuticals, Inc.
SO PCT Int. Appl., 182 pp.
CODEN: PIXXD2
DT Patent
LA English
FAN.CNT 1

	PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
PI	WO 2006044497	A2	20060427	WO 2005-US36752	20051012
	WO 2006044497	A3	20060908		
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	RW: AT, BE, BG, CH, CY, CZ, DE, DK, EE, ES, FI, FR, GB, GR, HU, IE, IS, IT, LT, LU, LV, MC, NL, PL, PT, RO, SE, SI, SK, TR, BF, BJ, CF, CG, CI, CM, GA, GN, GQ, GW, ML, MR, NE, SN, TD, TG, BW, GH, GM, KE, LS, MW, MZ, NA, SD, SL, SZ, TZ, UG, ZM, ZW, AM, AZ, BY, KG, KZ, MD, RU, TJ, TM				
	AU 2005295814	A1	20060427	AU 2005-295814	20051012
	CA 2583342	A1	20060427	CA 2005-2583342	20051012
	EP 1804794	A2	20070711	EP 2005-812233	20051012
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	CN 101068545	A	20071107	CN 2005-80034396	20051012
	US 20070197571	A1	20070823	US 2007-663388	20070321
	IN 2007CN01217	A	20070831	IN 2007-CN1217	20070323
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	WO 2005-US36752	W	20051012		
OS	MARPAT 144:432808				

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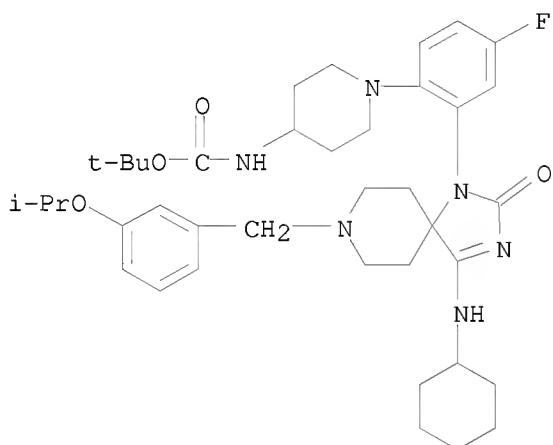
L4 ANSWER 1 OF 1 CAPLUS COPYRIGHT 2008 ACS on STN
IT 885115-63-9P 885115-64-0P
RL: RCT (Reactant); SPN (Synthetic preparation); PREP (Preparation); RACT
(Reactant or reagent)
(preparation of spiropiperidine compds. as β -secretase inhibitors for
treating Alzheimer's disease)
RN 885115-63-9 CAPLUS
CN Carbamic acid, [1-[4-fluoro-2-[8-[[3-(1-methylethoxy)phenyl]methyl]-2-oxo-

4-thioxo-1,3,8-triazaspiro[4.5]dec-1-ylphenyl]-4-piperidinyl-,
1,1-dimethylethyl ester (9CI) (CA INDEX NAME)



RN 885115-64-0 CAPLUS

CN Carbamic acid, [1-[2-[4-(cyclohexylamino)-8-[[3-(1-methylethoxy)phenyl]methyl]-2-oxo-1,3,8-triazaspiro[4.5]dec-3-en-1-yl]-4-fluorophenyl]-4-piperidinyl-, 1,1-dimethylethyl ester (9CI) (CA INDEX NAME)



=> file reg
COST IN U.S. DOLLARS
FULL ESTIMATED COST

SINCE FILE ENTRY	TOTAL SESSION
7.11	186.14

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10/589515

STRUCTURE FILE UPDATES: 28 APR 2008 HIGHEST RN 1017984-01-8
DICTIONARY FILE UPDATES: 28 APR 2008 HIGHEST RN 1017984-01-8

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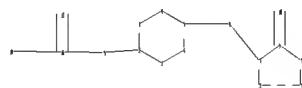
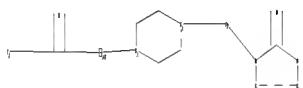
TSCA INFORMATION NOW CURRENT THROUGH January 9, 2008.

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<http://www.cas.org/support/stngen/stndoc/properties.html>

=>
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chain nodes :

8 14 15 16 17 19

ring nodes :

1 2 3 4 5 6 9 10 11 12 13

chain bonds :

2-15 5-8 8-13 9-14 15-16 16-17 16-19

ring bonds :

1-2 1-6 2-3 3-4 4-5 5-6 9-10 9-13 10-11 11-12 12-13

exact/norm bonds :

1-2 1-6 2-3 2-15 3-4 4-5 5-6 5-8 8-13 9-10 9-13 9-14 10-11 11-12
12-13 15-16 16-17 16-19

10/589515

G1:C,N

G2:O,N

Match level :

1:Atom 2:Atom 3:Atom 4:Atom 5:Atom 6:Atom 8:Atom 9:Atom 10:Atom 11:Atom
12:Atom 13:Atom 14:CLASS 15:CLASS 16:CLASS 17:CLASS 19:CLASS

Generic attributes :

8:

Saturation : Unsaturated
Number of Carbon Atoms : less than 7
Type of Ring System : Monocyclic

L5 STRUCTURE UPLOADED

=> s 15
SAMPLE SEARCH INITIATED 13:26:24 FILE 'REGISTRY'
SAMPLE SCREEN SEARCH COMPLETED - 18465 TO ITERATE

10.8% PROCESSED 2000 ITERATIONS 3 ANSWERS
INCOMPLETE SEARCH (SYSTEM LIMIT EXCEEDED)
SEARCH TIME: 00.00.01

FULL FILE PROJECTIONS: ONLINE **COMPLETE**
BATCH **COMPLETE**
PROJECTED ITERATIONS: 361163 TO 377437
PROJECTED ANSWERS: 238 TO 868

L6 3 SEA SSS SAM L5

=> s 15 sss full
FULL SEARCH INITIATED 13:26:31 FILE 'REGISTRY'
FULL SCREEN SEARCH COMPLETED - 367795 TO ITERATE

100.0% PROCESSED 367795 ITERATIONS 414 ANSWERS
SEARCH TIME: 00.00.04

L7 414 SEA SSS FUL L5

=> file caplus
COST IN U.S. DOLLARS SINCE FILE TOTAL
SESSION
FULL ESTIMATED COST 178.36 364.50

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=> s 17
 L8 1 L7

=> d 18 bib abs fhitstr

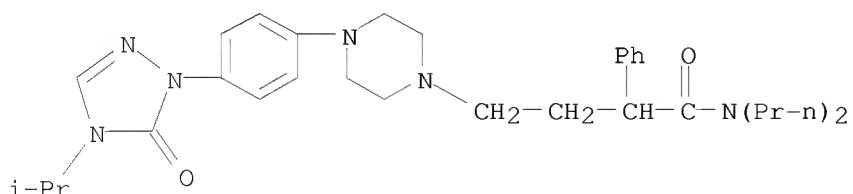
L8 ANSWER 1 OF 1 CAPLUS COPYRIGHT 2008 ACS on STN
 AN 2005:1004731 CAPLUS
 DN 143:306337
 TI Preparation of Mtp inhibiting aryl piperidines or piperazines substituted with 5-membered heterocycles
 IN Meerpoel, Lieven; Backx, Leo Jacobus Jozef; Jaroskova, Libuse; Roevens, Peter Walter Maria; Van, Der Veken Louis Jozef Elisabeth; Viellevoye, Marcel; Linders, Joannes Theodorus Maria
 PA Janssen Pharmaceutica N.V., Belg.; Meerpoel, Lieven; Backx, Leo Jacobus Jozef; Jaroskova, Libuse; Roevens, Peter Walter Maria; Van der Veken, Louis Jozef Elisabeth; Viellevoye, Marcel; Linders, Joannes Theodorus Maria
 SO PCT Int. Appl., 120 pp.
 CODEN: PIXXD2
 DT Patent
 LA English
 FAN.CNT 1

	PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
PI	WO 2005085226	A1	20050915	WO 2005-EP51010	20050307
	W: AE, AG, AL, AM, AT, AU, AZ, BA, BB, BG, BR, BW, BY, BZ, CA, CH, CN, CO, CR, CU, CZ, DE, DK, DM, DZ, EC, EE, EG, ES, FI, GB, GD, GE, GH, GM, HR, HU, ID, IL, IN, IS, JP, KE, KG, KP, KR, KZ, LC, LK, LR, LS, LT, LU, LV, MA, MD, MG, MK, MN, MW, MX, MZ, NA, NI, NO, NZ, OM, PG, PH, PL, PT, RO, RU, SC, SD, SE, SG, SK, SL, SM, SY, TJ, TM, TN, TR, TT, TZ, UA, UG, US, UZ, VC, VN, YU, ZA, ZM, ZW RW: BW, GH, GM, KE, LS, MW, MZ, NA, SD, SL, SZ, TZ, UG, ZM, ZW, AM, AZ, BY, KG, KZ, MD, RU, TJ, TM, AT, BE, BG, CH, CY, CZ, DE, DK, EE, ES, FI, FR, GB, GR, HU, IE, IS, IT, LT, LU, MC, NL, PL, PT, RO, SE, SI, SK, TR, BF, BJ, CF, CG, CI, CM, GA, GN, GQ, GW, ML, MR, NE, SN, TD, TG				
	AU 2005219580	A1	20050915	AU 2005-219580	20050307
	CA 2558506	A1	20050915	CA 2005-2558506	20050307
	EP 1751131	A1	20070214	EP 2005-716940	20050307
	R: AT, BE, BG, CH, CY, CZ, DE, DK, EE, ES, FI, FR, GB, GR, HU, IE, IS, IT, LI, LT, LU, MC, NL, PL, PT, RO, SE, SI, SK, TR, AL, BA,				

HR, LV, MK, YU				
CN 1930143	A	20070314	CN 2005-80007466	20050307
BR 2005008514	A	20070814	BR 2005-8514	20050307
JP 2007527897	T	20071004	JP 2007-502335	20050307
KR 2007014121	A	20070131	KR 2006-716095	20060810
US 20070191383	A1	20070816	US 2006-589515	20060816
MX 2006PA10253	A	20070216	MX 2006-PA10253	20060908
IN 2006DN05167	A	20070803	IN 2006-DN5167	20060908
NO 2006004601	A	20061010	NO 2006-4601	20061010
PRAI EP 2004-75771	A	20040310		
US 2004-556336P	P	20040325		
WO 2005-EP51010	W	20050307		
OS MARPAT 143:306337				
GI				

* STRUCTURE DIAGRAM TOO LARGE FOR DISPLAY - AVAILABLE VIA OFFLINE PRINT *

- AB Title compds. I [Z = (un)substituted -N=CH=, -CH=N-, -CH₂CH₂-, or -CH=CH-; X₁ = C or N; at least one of X₂ or X₃ = N and the other equals CH or C when the dotted line represents a bond; or both X₂ and X₃ = N; R₁ = (un)substituted alkyl, aryl, cycloalkyl, etc.; R₂ = H, halo, alkyl; A = divalent carbon bridge or when X₃ = CH, A may = NH which may be optionally substituted; B = (un)substituted amine, heterocycle, alkoxy, etc.], and their pharmaceutically acceptable salts, are prepared and disclosed as having apo β secretion/MTP inhibiting activity and concomitant lipid lowering activity. Thus, e.g., II was prepared by reaction of III (preparation given) with Et α -phenylacrylate. Selected compds. of the invention possessed pIC₅₀ values of 5.5-6.5 in assays studying inhibition of apo β secretion. The invention further relates to methods for preparing pharmaceutical compns. comprising said compds. as well as the use of said compds. as a medicine for the treatment of hyperlipidemia, obesity and type II diabetes (Formula (I)).
- IT 864928-10-9P
 RL: PAC (Pharmacological activity); SPN (Synthetic preparation); THU (Therapeutic use); BIOL (Biological study); PREP (Preparation); USES (Uses)
 (drug candidate; Preparation of Mtp inhibiting aryl piperidines or piperazines substituted with 5-membered heterocycles)
- RN 864928-10-9 CAPLUS
- CN 1-Piperazinebutanamide, 4-[4-[4,5-dihydro-4-(1-methylethyl)-5-oxo-1H-1,2,4-triazol-1-yl]phenyl]- α -phenyl-N,N-dipropyl- (CA INDEX NAME)



RE.CNT 2 THERE ARE 2 CITED REFERENCES AVAILABLE FOR THIS RECORD
 ALL CITATIONS AVAILABLE IN THE RE FORMAT

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COST IN U.S. DOLLARS	SINCE FILE ENTRY	TOTAL SESSION	
FULL ESTIMATED COST	5.93	370.43	
DISCOUNT AMOUNTS (FOR QUALIFYING ACCOUNTS)	SINCE FILE ENTRY	TOTAL SESSION	
CA SUBSCRIBER PRICE	-0.80	-0.80	

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FILE LAST UPDATED: 01 May 1997 (19970501/UP)

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L1 STRUCTURE uploaded
L2 0 S L1
L3 2 S L1 SSS FULL

FILE 'CAPLUS' ENTERED AT 13:21:33 ON 29 APR 2008

L4 1 S L3

FILE 'REGISTRY' ENTERED AT 13:25:58 ON 29 APR 2008

L5 STRUCTURE uploaded
L6 3 S L5
L7 414 S L5 SSS FULL

FILE 'CAPLUS' ENTERED AT 13:26:46 ON 29 APR 2008

L8 1 S L7

FILE 'CAOLD' ENTERED AT 13:27:15 ON 29 APR 2008

=> s 13
L9 0 L3

10/589515

=> s 17
L10 0 L7

=> file chemcats
COST IN U.S. DOLLARS SINCE FILE TOTAL
ENTRY SESSION
FULL ESTIMATED COST 0.46 370.89

DISCOUNT AMOUNTS (FOR QUALIFYING ACCOUNTS) SINCE FILE TOTAL
ENTRY SESSION
CA SUBSCRIBER PRICE 0.00 -0.80

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=> s 13 or 17
0 L3
0 L7
L11 0 L3 OR L7

=> log h
COST IN U.S. DOLLARS SINCE FILE TOTAL
ENTRY SESSION
FULL ESTIMATED COST 0.94 371.83

DISCOUNT AMOUNTS (FOR QUALIFYING ACCOUNTS) SINCE FILE TOTAL
ENTRY SESSION
CA SUBSCRIBER PRICE 0.00 -0.80

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STN INTERNATIONAL SESSION SUSPENDED AT 13:27:52 ON 29 APR 2008